Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101415\
Data File : BM002901.D
Acq On : 14 Oct 2015 16:38
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 15 01:02:13 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 14 17:36:42 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	-0.01
2	S	2-Fluorophenol	0.963	0.865	10.2	87	0.02
3	S	Phenol-d6	1.373	1.175	14.4	83	0.02
4		2-Chlorophenol	1.272	1.131	11.1	83	0.00
5	C	Phenol	1.349	1.203	10.8	83	0.00
6		2-Methylphenol	1.032	0.888	14.0	82	0.00
7		3+4-Methylphenols	1.277	1.112	12.9	81	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
9	S	Nitrobenzene-d5	0.276	0.253	8.3	83	0.00
10	C	2-Nitrophenol	0.126	0.105	16.7	78	0.00
11		2,4-Dimethylphenol	0.276	0.243	12.0	81	0.00
12	C	2,4-Dichlorophenol	0.215	0.190	11.6	77	0.02
13	C	4-Chloro-3-methylphenol	0.269	0.228	15.2	79	0.00
14	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
15	S	2,4,6-Tribromophenol	0.183	0.154	15.8	78	0.00
16	C	2,4,6-Trichlorophenol	0.326	0.268	17.8	75	0.00
17		2,4,5-Trichlorophenol	0.347	0.297	14.4	77	0.00
18	S	2-Fluorobiphenyl	1.636	1.522	7.0	83	0.00
19	P	2,4-Dinitrophenol	0.027	0.008#	70.4#	60	0.02
20	P	4-Nitrophenol	0.132	0.082	37.9#	66	0.00
21	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
22		4,6-Dinitro-2-methylphenol	0.025	0.025	0.0	76	0.00
23	C	Pentachlorophenol	0.038	0.023	39.5#	71	0.00
24	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
25	S	Terphenyl-d14	0.760	0.686	9.7	87	0.00
26	I	Perylene-d12	1.000	1.000	0.0	93	0.00

(#) = Out of Range SPCC's out = 1 CCC's out = 1